

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE110819\
 Data File : BE100715.D
 Acq On : 8 Nov 2019 13:16
 Operator : JU
 Sample : K5725-04DL 25X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SO-4DL

Manual Integrations
 APPROVED

mohammad
 11/11/2019 8:09:49 AM

Quant Time: Nov 08 14:40:48 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 23 16:49:33 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	2003	0.40	ng	0.00
7) Naphthalene-d8	10.64	136	9706	0.40	ng	0.01
13) Acenaphthene-d10	14.49	164	7522	0.40	ng	0.01
19) Phenanthrene-d10	17.20	188	14798	0.40	ng	0.00
27) Chrysene-d12	21.37	240	16010	0.40	ng	0.01
34) Perylene-d12	23.85	264	17955	0.40	ng	0.02

System Monitoring Compounds						
4) 2-Fluorophenol	5.45	112	92	0.01	ng	0.01
5) Phenol-d6	7.03	99	167	0.02	ng	0.02
8) Nitrobenzene-d5	9.00	82	1052	0.17	ng	0.01
11) 2-Methylnaphthalene-d10	12.22	152	512	0.03	ng	0.00
14) 2,4,6-Tribromophenol	15.98	330	10	0.01	ng	0.04
15) 2-Fluorobiphenyl	13.10	172	195	0.01	ng	0.00
25) Fluoranthene-d10	19.22	212	2699	0.01	ng	0.00
29) Terphenyl-d14	19.82	244	832	0.02	ng	0.00

Target Compounds						Qvalue
9) Naphthalene	10.69	128	69553	0.681	ng	# 98
12) 2-Methylnaphthalene	12.30	142	20821	1.197	ng	# 87
16) Acenaphthylene	14.20	152	16116	0.498	ng	# 74
17) Acenaphthene	14.54	154	8613	0.412	ng	# 74
18) Fluorene	15.51	166	13999	0.504	ng	# 62
23) Phenanthrene	17.24	178	119073	2.925	ng	# 94
24) Anthracene	17.34	178	28650	0.759	ng	# 80
26) Fluoranthene	19.26	202	239844	4.515	ng	# 99
28) Pyrene	19.61	202	247832	5.541	ng	# 94
30) Benzo(a)anthracene	21.35	228	156176	2.990	ng	# 95
31) Chrysene	21.40	228	119905	2.297	ng	# 98
32) Bis(2-ethylhexyl)phthalate	21.28	149	20807	0.190	ng	# 100
33) Indeno(1,2,3-cd)pyrene	26.47	276	131127	2.044	ng	# 96
35) Benzo(b)fluoranthene	23.09	252	215787	4.122	ng	# 100
36) Benzo(k)fluoranthene	23.13	252	85108m	1.581	ng	# 98
37) Benzo(a)pyrene	23.73	252	156212	3.205	ng	# 92
38) Dibenzo(a,h)anthracene	26.48	278	30870	0.613	ng	# 99
39) Benzo(a,h,i)perylene	27.28	276	136979	2.704	ng	# 99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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